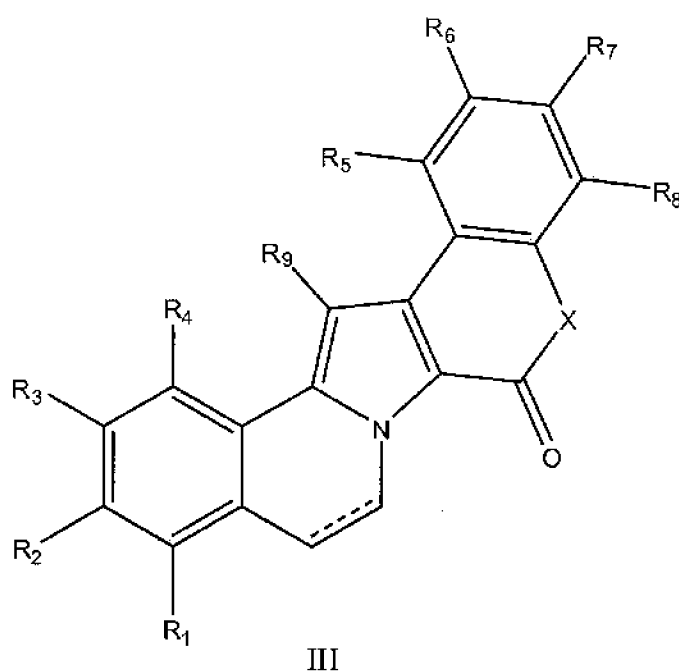


## AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions, and listings, of claims in the application:

**Listing of Claims:**

1. (currently amended) A compound of the general formula **III**:



wherein X is selected from the group consisting of NH, O and S;

wherein R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, and R<sub>8</sub> are each independently selected from the group consisting of H, OH, OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, C(=O)R', CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> alkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> haloalkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or

unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted aralkyl and substituted or unsubstituted heteroaromatic;

wherein R<sub>9</sub> is independently selected from the group consisting of H, OH, OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, C(=O)R', CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> alkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> haloalkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aralkyl, substituted or unsubstituted heteroaromatic, ~~and halogen~~ bromine, and iodine;

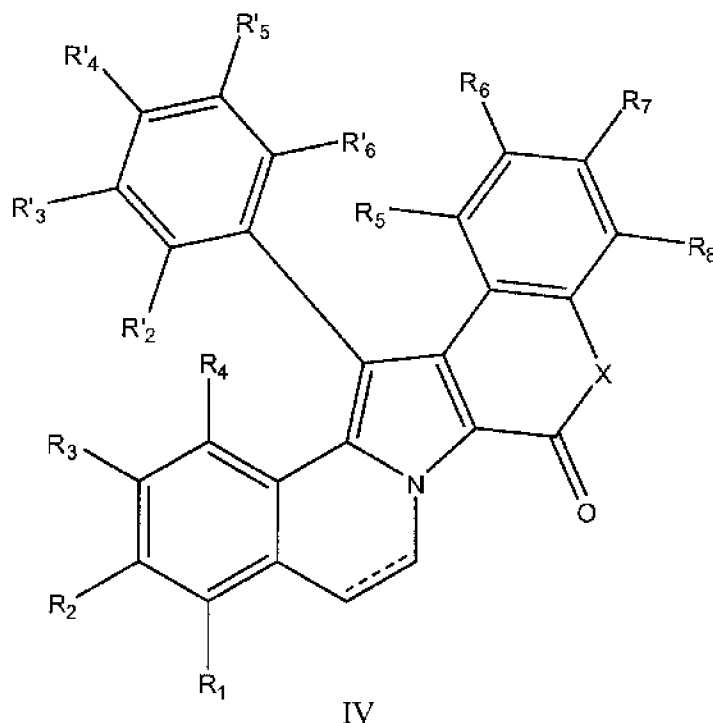
wherein each of the R' groups is independently selected from the group consisting of H, OH, NO<sub>2</sub>, NH<sub>2</sub>, SH, CN, halogen, C(=O)H, C(=O)CH<sub>3</sub>, CO<sub>2</sub>H, C(=O)R', substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>18</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>18</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkoxy, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> aminoalkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> aminoacid or aminoacids chain, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> thioalkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkylsulfinyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkylsulfonyl;

wherein the pairs of groups R<sub>1</sub> and R<sub>2</sub>, R<sub>2</sub> and R<sub>3</sub>, R<sub>3</sub> and R<sub>4</sub>, R<sub>3</sub> and R<sub>9</sub>, R<sub>4</sub> and R<sub>9</sub>, R<sub>9</sub> and R<sub>5</sub>, R<sub>9</sub> and R<sub>6</sub>, or R<sub>6</sub> and R<sub>7</sub>, R<sub>7</sub> and R<sub>8</sub> may be joined into a carbocyclic or heterocyclic ring system;

and the dotted line represents a single or double bond;

or a pharmaceutically acceptable salt or stereoisomer thereof.

2. (previously presented) A compound of the general formula IV:



wherein X is selected from the group consisting of NH, O and S;

wherein R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>8</sub>, R'<sub>2</sub>, R'<sub>3</sub>, R'<sub>4</sub>, R'<sub>5</sub>, and R'<sub>6</sub> are each independently selected from the group consisting of H, OH, OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, C(=O)R', CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> alkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> haloalkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted aralkyl and substituted or unsubstituted heteroaromatic;

wherein  $R_7$  is independently selected from the group consisting of  $OR'$ ,  $SH$ ,  $SR'$ ,  $SOR'$ ,  $SO_2R'$ ,  $NHR'$ ,  $N(R')_2$ ,  $N=R'$ ,  $NHCOR'$ ,  $N(COR')_2$ ,  $NHSO_2R'$ ,  $NO_2$ ,  $PO(R')_2$ ,  $PO_2R'$ ,  $C(=O)H$ ,  $C(=O)R'$ ,  $CO_2H$ ,  $CO_2R'$ ,  $OPO(R')_2$ ,  $OPO_2R'$ ,  $OC(=O)H$ ,  $OC(=O)R'$ ,  $N=C(R')_2$ , substituted or unsubstituted  $C_1-C_{12}$  alkyl, substituted or unsubstituted  $C_1-C_{12}$  haloalkyl, substituted or unsubstituted  $C_2-C_{12}$  alkenyl, substituted or unsubstituted  $C_2-C_{12}$  alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted aralkyl and substituted or unsubstituted heteroaromatic;

with the proviso that  $R_7$  is not Me, Et, Pr, COMe, OH, OMe, OAc, O<sup>i</sup>Pr or OBn when X is O;

wherein each of the  $R'$  groups is independently selected from the group consisting of H, OH,  $NO_2$ ,  $NH_2$ ,  $SH$ ,  $CN$ , halogen,  $C(=O)H$ ,  $C(=O)R'$ ,  $C(=O)CH_3$ ,  $CO_2H$ , substituted or unsubstituted  $C_1-C_{18}$  alkyl, substituted or unsubstituted  $C_2-C_{18}$  alkenyl, substituted or unsubstituted  $C_2-C_{18}$  alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted  $C_1-C_{18}$  alkoxy, substituted or unsubstituted  $C_1-C_{18}$  aminoalkyl, substituted or unsubstituted  $C_1-C_{18}$  aminoacid or aminoacids chain, substituted or unsubstituted  $C_1-C_{18}$  thioalkyl, substituted or unsubstituted  $C_1-C_{18}$  alkylsulfinyl, substituted or unsubstituted  $C_1-C_{18}$  alkylsulfonyl;

wherein the pairs of groups  $R_1$  and  $R_2$ ,  $R_2$  and  $R_3$ ,  $R_3$  and  $R_4$ ,  $R_6$  and  $R_7$ , or  $R_7$  and  $R_8$  may be joined into a carbocyclic or heterocyclic ring system;

and the dotted line represents a single or double bond;

or a pharmaceutically acceptable salt or stereoisomer thereof.

3. (previously presented) A compound according to claim 1 or 2 wherein X is O or NH.

4. (previously presented) A compound according to claim 1 or 2 wherein X is O.
5. (previously presented) A compound according to claim 1 or 2 wherein the dotted line is a double bond.
6. (previously presented) A compound according to claim 1 wherein each of  $R_1$ - $R_8$  is independently selected from H,  $OR'$ , and  $OC(=O)R'$ .
7. (previously presented) A compound according to claim 1 or 2 wherein  $R_3$  is selected from the group consisting of H, OH, and  $OR'$ , with the proviso that when  $R_3$  is  $OR'$ , then  $R'$  is selected from a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl.
8. (previously presented) A compound according to claim 1 or 2 wherein  $R_4$ ,  $R_5$ ,  $R_6$  and  $R_8$  are each independently selected from the group consisting of H and  $OR'$ , with the proviso that when  $R_4$ ,  $R_5$ ,  $R_6$  or  $R_8$  is  $OR'$ , then  $R'$  is selected from a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl.
9. (previously presented) A compound according to claim 8 wherein  $R_4$ ,  $R_5$  and  $R_8$  are H.
10. (previously presented) A compound according to claim 1 wherein  $R_1$ ,  $R_2$  and  $R_7$  are each independently selected from the group consisting of H, OH,  $OR'$ ,  $OC(=O)R'$ ,  $SO_2R'$ ,  $PO(R')_2$ , substituted or unsubstituted  $C_1$ - $C_{12}$  alkyl,  $NO_2$ , and  $NH_2$ , with the proviso that when  $R_1$ ,  $R_2$  or  $R_7$  are  $OR'$ , then  $R'$  is selected from a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl.

11. (previously presented) A compound according to claim 10 wherein  $R_1$ ,  $R_2$  and  $R_7$  are  $OC(=O)R'$  wherein  $R'$  is a substituted or unsubstituted aminoacid or aminoacids chain.

12. (previously presented) A compound according to claim 2 wherein  $R'_2$ ,  $R'_3$  and  $R'_6$  are each independently selected from the group consisting of H and  $OR'$ , wherein  $R'$  is a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl.

13. (previously presented) A compound according to claim 2 wherein  $R'_5$  is selected from the group consisting of H and  $OR'$ , wherein  $R'$  is a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl.

14. (previously presented) A compound according to claim 2 wherein  $R'_4$  is selected from the group consisting of H, OH,  $OR'$ ,  $OC(=O)R'$ ,  $SO_2R'$ ,  $PO(R')_2$ , substituted or unsubstituted  $C_1$ - $C_{12}$  alkyl,  $NO_2$ , and  $NH_2$ , with the proviso that when  $R'_4$  is  $OR'$ , then  $R'$  is selected from a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl.

15. (currently amended) A compound according to claim 14 ~~wherein  $R'_4$  is  $OC(=O)R'$  wherein  $R'_4$  is  $OR'$  and~~ wherein  $R'$  is a substituted or unsubstituted aminoacid or aminoacids chain.

16. (previously presented) A compound according to claim 1 or 2 wherein at least one of  $R_1$ - $R_8$  and  $R'_2$ - $R'_6$  is not H, OH,  $OCH_3$ , and  $SO_3Na$ .

17. (previously presented) A pharmaceutical composition comprising a compound as defined in

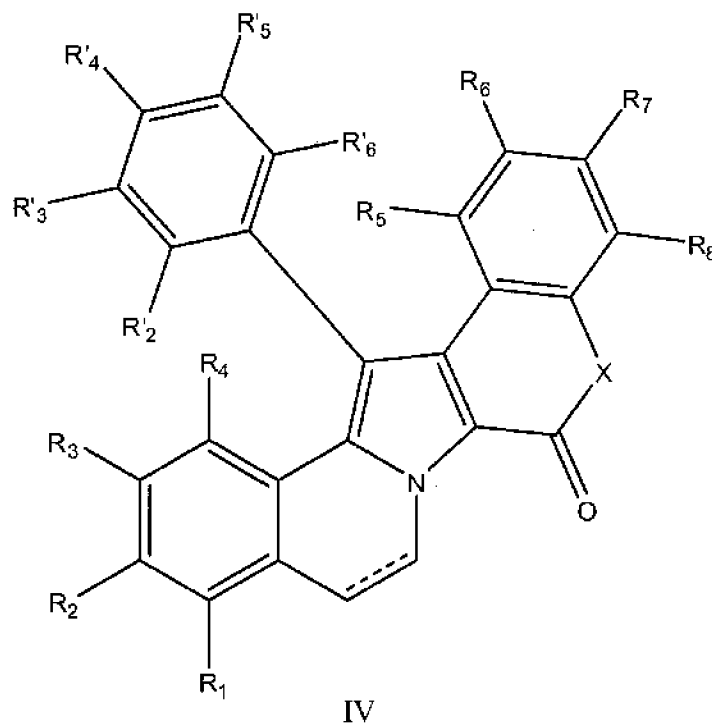
claim 1 or 2 or a pharmaceutically acceptable salt or stereoisomer thereof, and a pharmaceutically acceptable diluent or carrier.

18. (canceled)

19. (previously presented) A method of treating a tumor which comprises administering to a human an effective amount of a compound as defined in claim 1 or 2 or a pharmaceutically acceptable salt or stereoisomer thereof.

20. (previously presented) A method of inhibiting topoisomerase I comprising administering to a human an amount effective for inhibiting topoisomerase I of a compound as defined in claim 1 or 2 or a pharmaceutically acceptable salt or stereoisomer thereof.

21. (previously presented) A compound of the general formula **IV**:



wherein X is selected from the group consisting of NH, O and S;

wherein  $R_1$ ,  $R_2$ ,  $R_3$ ,  $R_4$ ,  $R_5$ ,  $R_6$ ,  $R_8$ ,  $R'_2$ ,  $R'_3$ ,  $R'_4$ ,  $R'_5$ , and  $R'_6$  are each independently selected from the group consisting of H, OH, OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, C(=O)R', CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> alkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> haloalkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted aralkyl and substituted or unsubstituted heteroaromatic;

wherein  $R_7$  is independently selected from the group consisting of OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub>



haloalkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted alkyl and substituted or unsubstituted heteroaromatic;

with the proviso that R<sub>7</sub> is not OH, OMe, OAc, O<sup>i</sup>Pr or OBn when X is O;

wherein each of the R' groups is independently selected from the group consisting of H, OH, NO<sub>2</sub>, NH<sub>2</sub>, SH, CN, halogen, C(=O)H, C(=O)CH<sub>3</sub>, CO<sub>2</sub>H, C(=O)R', substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>18</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>18</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkoxy, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> aminoalkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> aminoacid or aminoacids chain, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> thioalkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkylsulfinyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkylsulfonyl;

wherein the pairs of groups R<sub>1</sub> and R<sub>2</sub>, R<sub>2</sub> and R<sub>3</sub>, R<sub>3</sub> and R<sub>4</sub>, R<sub>6</sub> and R<sub>7</sub>, or R<sub>7</sub> and R<sub>8</sub> may be joined into a carbocyclic or heterocyclic ring system;

and the dotted line represents a single or double bond;

or a pharmaceutically acceptable salt or stereoisomer thereof.

22. (previously presented) A compound according to claim 7 wherein R<sub>3</sub> is methoxy.

23. (previously presented) A compound according to claim 13 wherein R'<sub>5</sub> is methoxy.

24. (previously presented) A compound according to claim 16 wherein at least two of  $R_1$ - $R_8$  and  $R'_2$ - $R'_6$  are not H, OH,  $OCH_3$ , or  $SO_3Na$ .

25. (previously presented) A compound according to claim 11 wherein  $R'$  is an aminoacid or aminoacids chain substituted with a cationic group.

26. (previously presented) A compound according to claim 15 wherein  $R'$  is an aminoacid or aminoacids chain substituted with a cationic group.

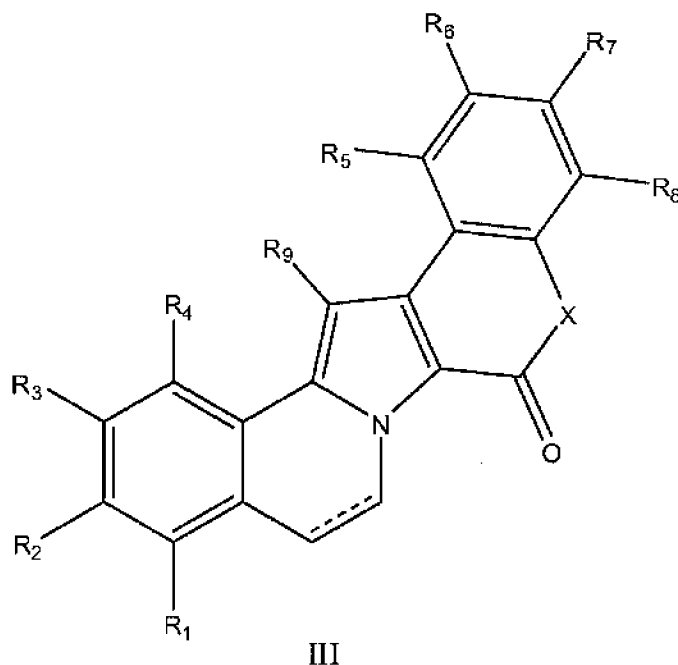
27. (previously presented) A compound according to claim 2 or 21 wherein each of  $R_1$ - $R_6$  and  $R_8$  is independently selected from H,  $OR'$ , and  $OC(=O)R'$  and wherein  $R_7$  is selected from  $OR'$  and  $OC(=O)R'$ .

28. (previously presented) A compound according to claim 2 or 21 wherein  $R_1$  and  $R_2$  are each independently selected from the group consisting of H, OH,  $OR'$ ,  $OC(=O)R'$ ,  $SO_2R'$ ,  $PO(R')_2$ , substituted or unsubstituted  $C_1$ - $C_{12}$  alkyl,  $NO_2$ , and  $NH_2$ , with the proviso that when  $R_1$  or  $R_2$  are  $OR'$ , then  $R'$  is selected from a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl; and

wherein  $R_7$  is selected from the group consisting of  $OR'$ ,  $OC(=O)R'$ ,  $SO_2R'$ ,  $PO(R')_2$ , substituted or unsubstituted  $C_1$ - $C_{12}$  alkyl,  $NO_2$ , and  $NH_2$ , with the proviso that when  $R_7$  is  $OR'$ , then  $R'$  is selected from a substituted or unsubstituted  $C_1$ - $C_{18}$  alkyl group.

29. (previously presented) A compound according to claim 12 wherein  $R'_2$ ,  $R'_3$ , and  $R'_6$  are H.

30. (new) A compound of the general formula **III**:



wherein X is selected from the group consisting of NH, O and S;

wherein R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, and R<sub>8</sub> are each independently selected from the group consisting of H, OH, OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, C(=O)R', CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> alkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> haloalkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted aralkyl and substituted or unsubstituted heteroaromatic;

wherein R<sub>9</sub> is independently selected from the group consisting of H, OH, OR', SH, SR', SOR', SO<sub>2</sub>R', NHR', N(R')<sub>2</sub>, N=R', NHCOR', N(COR')<sub>2</sub>, NHSO<sub>2</sub>R', NO<sub>2</sub>, PO(R')<sub>2</sub>, PO<sub>2</sub>R', C(=O)H, C(=O)R', CO<sub>2</sub>H, CO<sub>2</sub>R', OPO(R')<sub>2</sub>, OPO<sub>2</sub>R', OC(=O)H, OC(=O)R', N=C(R')<sub>2</sub>, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> alkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>12</sub> haloalkyl, substituted or

unsubstituted C<sub>2</sub>-C<sub>12</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>12</sub> alkynyl, substituted or unsubstituted aralkyl, substituted or unsubstituted heteroaromatic, and halogen;

wherein each of the R' groups is independently selected from the group consisting of H, OH, NO<sub>2</sub>, NH<sub>2</sub>, SH, CN, halogen, C(=O)H, C(=O)CH<sub>3</sub>, CO<sub>2</sub>H, C(=O)R', substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkyl, substituted or unsubstituted C<sub>2</sub>-C<sub>18</sub> alkenyl, substituted or unsubstituted C<sub>2</sub>-C<sub>18</sub> alkynyl, substituted or unsubstituted aryl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkoxy, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> aminoalkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> aminoacid or aminoacids chain, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> thioalkyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkylsulfinyl, substituted or unsubstituted C<sub>1</sub>-C<sub>18</sub> alkylsulfonyl;

wherein the pairs of groups R<sub>1</sub> and R<sub>2</sub>, R<sub>2</sub> and R<sub>3</sub>, R<sub>3</sub> and R<sub>4</sub>, R<sub>3</sub> and R<sub>9</sub>, R<sub>4</sub> and R<sub>9</sub>, R<sub>9</sub> and R<sub>5</sub>, R<sub>9</sub> and R<sub>6</sub>, or R<sub>6</sub> and R<sub>7</sub>, R<sub>7</sub> and R<sub>8</sub> may be joined into a carbocyclic or heterocyclic ring system;

and the dotted line represents a single or double bond;

or a pharmaceutically acceptable salt or stereoisomer thereof.